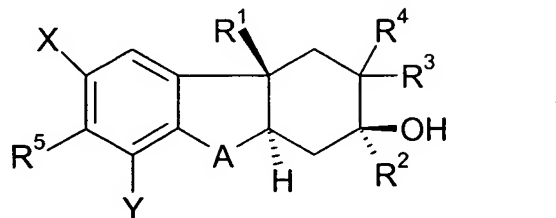


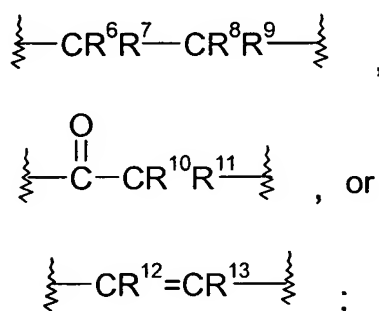
AMENDMENTS TO THE CLAIMS

This listing of claims will replace all prior versions and listings of claims in the application.

1. (Original) A compound of the formula



wherein A is of the formula



X and Y are each independently hydrogen, fluoro, chloro, bromo, or (C₁-C₆)alkyl;

R¹ is (C₂-C₆)alkyl, (C₃-C₆)alkenyl, or optionally substituted benzyl; wherein said benzyl may be optionally substituted with one to three substituents independently selected from HO-, (C₁-C₆)alkyl-O-, halo and amino;

R² is (C₁-C₆)alkyl, (C₂-C₆)alkenyl, (C₃-C₆)alkynyl, (C₃-C₁₀)cycloalkyl, (C₆-C₁₀)aryl, (C₁-C₉)heterocyclyl, (C₁-C₉)heteroaryl, (C₆-C₁₀)aryl(C₁-C₄)alkyl, (C₁-C₉)heterocyclyl-(C₁-C₄)alkyl, (C₁-C₉)heteroaryl-(C₁-C₄)alkyl, or (C₃-C₁₀)cycloalkyl-(C₁-C₄)alkyl; wherein each of the aforesaid groups may optionally be substituted with one to three substituents independently selected from halo, (C₁-C₆)alkyl, (C₁-C₆)alkoxy, or -CF₃;

R³ is hydrogen, (C₁-C₆)alkyl, (C₂-C₆)alkenyl, (C₂-C₆)alkynyl, (C₃-C₁₀)cycloalkyl, (C₁-C₉)heterocyclyl, (C₁-C₉)heteroaryl, or (C₆-C₁₀)aryl; wherein each of the aforesaid groups may be optionally substituted with one to three substituents independently selected from HO-, (C₁-C₆)alkyl-O-, halo and amino;

R⁴ is HO- or R¹⁴R¹⁵N-;

R⁵ is a radical selected from the group consisting of hydrogen, halo, (C₁-C₆)alkyl, (C₂-C₆)alkenyl, (C₃-C₆)alkynyl, (C₃-C₁₀)cycloalkyl, (C₆-C₁₀)aryl-, (C₁-C₉)heteroaryl-, (C₁-C₉)heterocyclic-, -OH, (C₁-C₆)alkyl-O-, (C₃-C₁₀)cycloalkyl-O-, (C₆-C₁₀)aryl-O-, (C₁-C₉)heteroaryl-O-, (C₁-C₉)heterocyclic-O-, (C₃-C₁₀)cycloalkyl-(C₁-C₆)alkyl-O-, (C₆-C₁₀)aryl-(C₁-C₆)alkyl-O-, (C₁-C₉)heteroaryl-(C₁-C₆)alkyl-O-, (C₁-C₉)heterocyclic-(C₁-C₆)alkyl-O-, R¹⁶R¹⁷N-(C=O)-, R¹⁶-(C=O)-(R²⁵-N)-, R¹⁶R¹⁷-N-SO₂-, R¹⁸-SO₂-, R¹⁸-SO₂-(NR¹⁹)-, R¹⁸-SO₃-, -C≡N, R¹⁸-(C=O)-O-, R¹⁸-(C=O)-, R¹⁶R¹⁷N-(C=O)-O-, R¹⁶R¹⁷N-(C=O)-(R²⁵-N)-, R¹⁹-O-(C=O)-(R²⁵-N)-, and R¹⁹-O-(C=O)-; wherein each of said (C₁-C₆)alkyl, (C₃-C₁₀)cycloalkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl, (C₁-C₉)heterocyclic moieties of said (C₁-C₆)alkyl, (C₆-C₁₀)aryl-, (C₁-C₉)heteroaryl-, (C₁-C₉)heterocyclic-, (C₁-C₆)alkyl-O-, (C₃-C₁₀)cycloalkyl-O-, (C₆-C₁₀)aryl-O-, (C₁-C₉)heteroaryl-O-, (C₁-C₉)heterocyclic-O-, (C₃-C₁₀)cycloalkyl-(C₁-C₆)alkyl-O-, (C₆-C₁₀)aryl-(C₁-C₆)alkyl-O-, (C₁-C₉)heteroaryl-(C₁-C₆)alkyl-O- and (C₁-C₉)heterocyclic-(C₁-C₆)alkyl-O- radicals, may optionally be substituted with one to three substituents independently selected from the group consisting of (C₁-C₆)alkyl, (C₂-C₆)alkenyl, (C₂-C₆)alkynyl, (C₃-C₁₀)cycloalkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl(CH₂)_n-, (C₁-C₉)heterocyclic, halo, HO-, HO-(C=O)-, R²⁰-O-(C=O)-, R²¹-(C=O)-, R²²-CO₂-, N≡C-, R²³R²⁴N-, R²³R²⁴N-(C₁-C₆)alkyl-, R²³R²⁴N-(C=O)-, R²³R²⁴-N-SO₂-, R²¹-SO₂-, R²¹-SO₂-(NR²¹)-, R²¹-SO₃-, R²¹(C=O)-NH-, R²¹(C=O)-[N-(C₁-C₆)alkyl]-; R²¹(C=O)-NH-(C₁-C₆)alkyl-; and R²¹(C=O)-[N-(C₁-C₆)alkyl]-(C₁-C₆)alkyl-; wherein said (C₃-C₁₀)cycloalkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl(CH₂)_n-, (C₁-C₉)heterocyclic substituents may optionally be substituted on a ring carbon or nitrogen by one to three members per ring independently selected from halo, (C₁-C₆)alkyl, and (C₁-C₆)alkoxy;

n is an integer from zero to four;

each of R⁶, R⁷, R⁸ and R⁹ is independently selected from the group consisting of hydrogen, (C₁-C₆)alkyl, fluoro and -OH;

each of R¹⁰ and R¹¹ is independently selected from the group consisting of hydrogen and (C₁-C₆)alkyl;

each of R¹² and R¹³ is independently selected from the group consisting of hydrogen, fluoro and (C₁-C₆)alkyl;

each of R¹⁴ and R¹⁵ is independently selected from hydrogen or (C₁-C₄)alkyl;

each of R¹⁶ and R¹⁷ is independently selected from hydrogen, (C₁-C₆)alkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl, (C₁-C₉)heterocyclic, (C₁-C₉)heteroaryl(C₁-C₆)alkyl, (C₆-C₁₀)aryl(C₁-C₆)alkyl, (C₁-C₉)heterocyclic(C₁-C₆)alkyl, HO-(C₁-C₆)alkyl, amino-(C₁-C₆)alkyl-, (C₁-C₆)alkylamino-(C₁-C₆)alkyl-, and [(C₁-C₆)alkyl]₂amino-(C₁-C₆)alkyl-; wherein said each of said (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl, and (C₁-C₉)heterocyclic moieties of said (C₆-C₁₀)aryl-, (C₁-C₉)heteroaryl-, (C₁-C₉)heterocyclic-, (C₆-C₁₀)aryl-(C₁-C₆)alkyl, (C₁-C₉)heteroaryl-(C₁-C₆)alkyl and (C₁-C₉)heterocyclic-(C₁-C₆)alkyl, may optionally be substituted with one to three substituents independently selected from the group consisting of halo, (C₁-C₆)alkyl or (C₁-C₆)alkoxy, or R¹⁶ and R¹⁷ are taken together to form an azetidiny, pyrrolidiny, piperidiny, piperaziny, (C₁-C₆)alkyl-piperaziny, or morpholiny ring;

R¹⁸ is hydrogen, (C₁-C₆)alkyl, (C₆-C₁₀)aryl or (C₁-C₉)heteroaryl; wherein said (C₁-C₆)alkyl may optionally be substituted with a substituent selected from the group consisting of HO-, amino, (C₁-C₆)alkylamino, [(C₁-C₆)alkyl]₂amino, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl, (C₁-C₉)heterocyclic, (C₁-C₆)alkoxy, HO-(C=O)-, (C₁-C₆)alkyl-O-(C=O)-, (C₁-C₆)alkyl-(C=O)-, N≡C-, [(C₁-C₆)alkyl]₂N-(C=O)- and (C₁-C₆)alkyl(C=O)-NH-;

R¹⁹ is hydrogen or (C₁-C₆)alkyl;

R²⁰ is hydrogen or (C₁-C₆)alkyl;

R²¹ is hydrogen or (C₁-C₆)alkyl;

R²² is hydrogen or (C₁-C₆)alkyl;

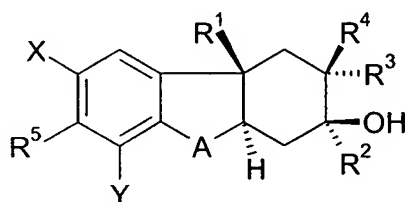
each of R²³ and R²⁴ is independently selected from hydrogen, (C₁-C₆)alkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl, (C₁-C₉)heterocyclic, (C₁-C₉)heteroaryl(C₁-C₆)alkyl, (C₆-C₁₀)aryl(C₁-C₆)alkyl, (C₁-C₉)heterocyclic(C₁-C₆)alkyl, HO-(C₁-C₆)alkyl, N≡C-(C₁-C₆)alkyl, amino-(C₁-C₆)alkyl-, (C₁-C₆)alkylamino-(C₁-C₆)alkyl-, and [(C₁-C₆)alkyl]₂amino-(C₁-C₆)alkyl-; wherein said each of said (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl, and (C₁-C₉)heterocyclic moieties of said (C₆-C₁₀)aryl-, (C₁-C₉)heteroaryl-, (C₁-C₉)heterocyclic-, (C₆-C₁₀)aryl-(C₁-C₆)alkyl, (C₁-C₉)heteroaryl-(C₁-C₆)alkyl and (C₁-C₉)heterocyclic-(C₁-C₆)alkyl, may optionally be substituted with one to three substituents independently selected from the group consisting of halo, (C₁-C₆)alkyl or (C₁-C₆)alkoxy, or R²³ and R²⁴ are taken together to

form an azetidiny, pyrrolidiny, piperidiny, piperaziny, (C₁-C₆)alkyl-piperaziny, or morpholiny ring;

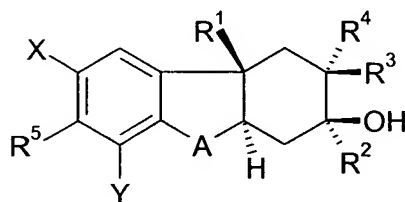
R²⁵ is hydrogen or (C₁-C₆)alkyl;

or a pharmaceutically acceptable salt thereof.

2. (Currently Amended)
compound has the formula

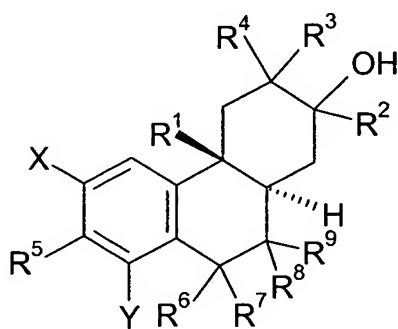


3. (Currently Amended)
compound has the formula



4. (Original)
formula

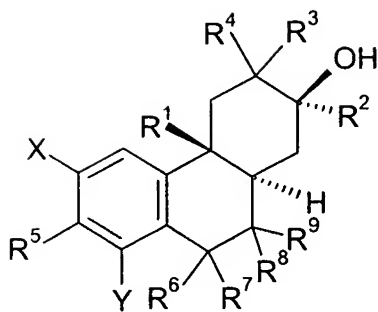
A compound according to claim 1, wherein said compound has the



1a

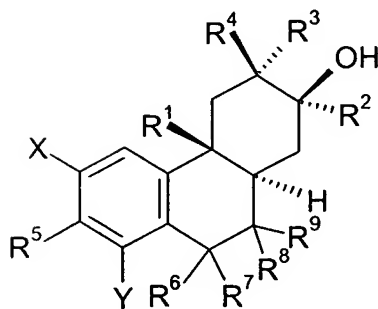
5. (Currently Amended)
compound has the formula

A compound according to claim 1, wherein said



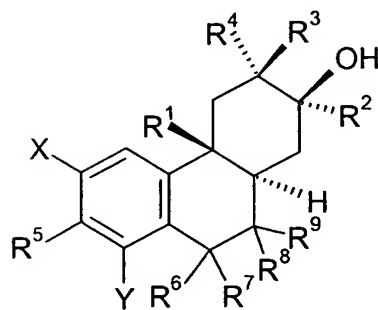
6. (Currently Amended)
compound has the formula

A compound according to claim 1, wherein said



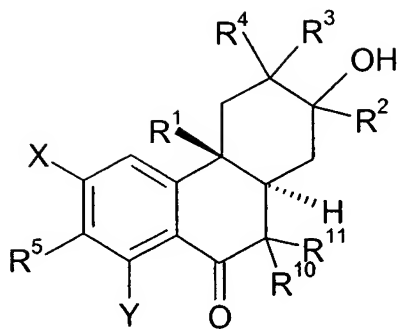
7. (Currently Amended)
compound has the formula

A compound according to claim 1, wherein said



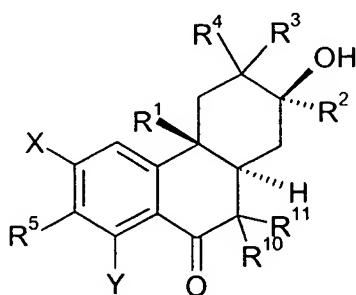
8. (Withdrawn)
has the formula

A compound according to claim 1, wherein said compound

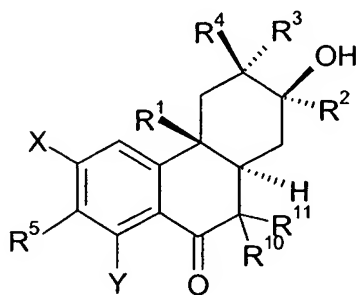


1b

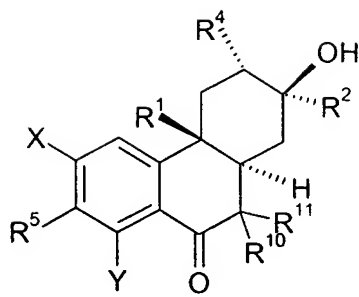
9. (Withdrawn-Currently Amended) A compound according to claim 1, wherein said compound has the formula



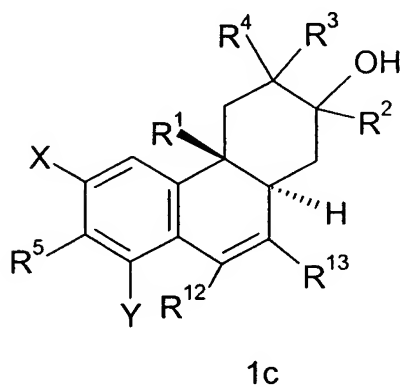
10. (Withdrawn-Currently Amended) A compound according to claim 1, wherein said compound has the formula



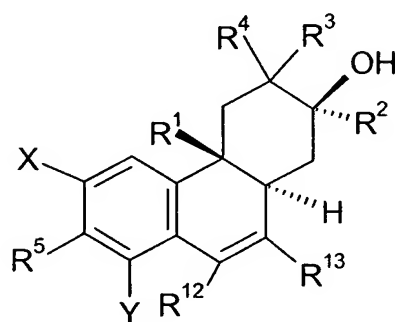
11. (Withdrawn-Currently Amended) A compound according to claim 1, wherein said compound has the formula



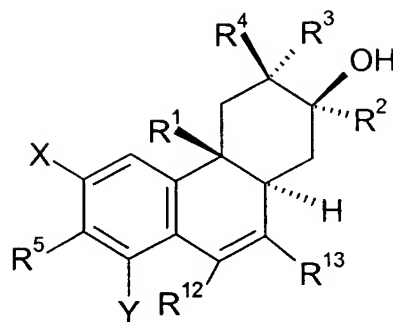
12. (Original) A compound according to claim 1, wherein said compound has the formula



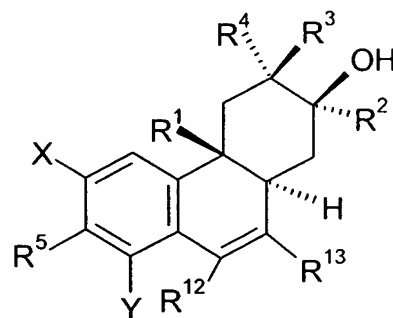
13. (Withdrawn-Currently Amended) A compound according to claim 1, wherein said compound has the formula



14. (Withdrawn-Currently Amended) A compound according to claim 1, wherein said compound has the formula



15. (Withdrawn-Currently Amended) A compound according to claim 1, wherein said compound has the formula



16. (Currently Amended) A compound according to claim 1 ~~any of the foregoing claims~~, wherein R¹ is ethyl or allyl.

17. (Currently Amended) A compound according to claim 1 ~~any of the foregoing claims~~, wherein R² is optionally substituted (C₆-C₁₀)aryl.

18. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is optionally substituted (C₁-C₉)heteroaryl.

19. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is optionally substituted (C₃-C₅)heteroaryl.

20. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is optionally substituted (C₁-C₉)heterocyclyl.

21. (Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is optionally substituted phenyl.
22. (Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is phenyl.
23. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is optionally substituted thiazolyl.
24. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is optionally substituted pyridyl.
25. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is optionally substituted oxazolyl.
26. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is optionally substituted pyridin-2-yl.
27. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is optionally substituted thiazol-2-yl.
28. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is optionally substituted oxazol-2-yl.
29. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is pyridin-2-yl; optionally substituted with a substituent selected from halo, CF₃, and (C₁-C₆)alkyl.

30. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is thiazol-2-yl; optionally substituted with a substituent selected from halo, CF₃, or (C₁-C₆)alkyl.
31. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is oxazol-2-yl; optionally substituted with a substituent selected from halo, CF₃, or (C₁-C₆)alkyl.
32. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is pyridin-2-yl.
33. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is thiazol-2-yl.
34. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is oxazol-2-yl.
35. (Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is (C₃-C₆)alkynyl.
36. (Currently Amended) A compound according to claim 1 ~~claims 1-16~~, wherein R² is (C₂-C₆)alkenyl.
37. (Currently Amended) A compound according to claim 1 ~~any of the foregoing claims~~, wherein R³ is hydrogen.
38. (Currently Amended) A compound according to claim 1 ~~claims 1-36~~, wherein R³ is (C₁-C₆)alkyl optionally substituted with a substituent selected from halo or hydroxy.
39. (Currently Amended) A compound according to claim 1 ~~claims 1-36~~, wherein R³ is methyl, ethyl or propyl.

40. (Currently Amended) A compound according to claim 1 ~~claims 1-36~~, wherein R³ is methyl.
41. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-36~~, wherein R³ is optionally substituted (C₁-C₉)heteroaryl.
42. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-36~~, wherein R³ is optionally substituted (C₁-C₉)heterocyclyl.
43. (Currently Amended) A compound according to claim 1 ~~claims 1-36~~, wherein R³ is optionally substituted (C₆-C₁₀)aryl.
44. (Currently Amended) A compound according to claims 1, 4, 5, 6, and 7 ~~any of the foregoing claims~~, wherein R⁴ is HO-.
45. (Currently Amended) A compound according to claim 1 ~~claims 1-36~~, wherein R⁴ is R¹⁴R¹⁵N-.
46. (Withdrawn-Currently Amended) A compound according to claim 1 ~~any of the foregoing claims~~, wherein R⁵ is -OH.
47. (Currently Amended) A compound according to claim 1 ~~claims 1-45~~, wherein R⁵ is (C₁-C₆)alkyl-O-, (C₃-C₁₀)cycloalkyl-O-, (C₆-C₁₀)aryl-O-, (C₁-C₉)heteroaryl-O-, or (C₁-C₉)heterocyclic-O-, wherein each of said (C₁-C₆)alkyl, (C₃-C₁₀)cycloalkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl, (C₁-C₉)heterocyclic moieties of said (C₁-C₆)alkyl-O-, (C₃-C₁₀)cycloalkyl-O-, (C₆-C₁₀)aryl-O-, (C₁-C₉)heteroaryl-O-, (C₁-C₉)heterocyclic-O- radicals may optionally be substituted with one to three substituents independently selected from (C₁-C₆)alkyl, (C₂-C₆)alkenyl, (C₂-C₆)alkynyl, (C₃-C₁₀)cycloalkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl, (C₁-C₉)heterocyclic, halo, HO-, HO-(C=O)-, R²¹-(C=O)-, R²²-CO₂-, N≡C-, R²³R²⁴N-, R²³R²⁴N-(C=O)-, R²¹(C=O)-NH-, R²¹(C=O)-[N-(C₁-C₆)alkyl]-.

48. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-45~~, wherein R⁵ is optionally substituted (C₆-C₁₀)aryl-, (C₁-C₉)heteroaryl-, (C₁-C₉)heterocyclic-, (C₆-C₁₀)aryl-(C₁-C₆)alkyl, (C₁-C₉)heteroaryl-(C₁-C₆)alkyl or (C₁-C₉)heterocyclic-(C₁-C₆)alkyl; optionally substituted with one to three substituents independently selected from (C₁-C₆)alkyl, (C₂-C₆)alkenyl, (C₂-C₆)alkynyl, (C₃-C₁₀)cycloalkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl, (C₁-C₉)heterocyclic, halo, HO-, HO-(C=O)-, R²¹-(C=O)-, R²²-CO₂-, N≡C-, R²³R²⁴N-, R²³R²⁴N-(C=O)-, R²¹(C=O)-NH-, R²¹(C=O)-[N-(C₁-C₆)alkyl]-.

49. (Currently Amended) A compound according to claim 1 ~~claims 1-45~~, wherein R⁵ is (C₆-C₁₀)aryl-(C₁-C₆)alkyl-O-, (C₁-C₉)heteroaryl-(C₁-C₆)alkyl-O-, (C₁-C₉)heterocyclic-(C₁-C₆)alkyl-O-, wherein each of said (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl, (C₁-C₉)heterocyclic moieties of said (C₆-C₁₀)aryl-(C₁-C₆)alkyl-O-, (C₁-C₉)heteroaryl-(C₁-C₆)alkyl-O-, and (C₁-C₉)heterocyclic-(C₁-C₆)alkyl-O-, may optionally be substituted with one to three substituents independently selected from the group consisting of (C₁-C₆)alkyl, (C₂-C₆)alkenyl, (C₂-C₆)alkynyl, (C₃-C₁₀)cycloalkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl(CH₂)_n-, (C₁-C₉)heterocyclic, halo, HO-, HO-(C=O)-, R²⁰-O-(C=O)-, R²¹-(C=O)-, R²²-CO₂-, N≡C-, R²³R²⁴N-, R²³R²⁴N-(C₁-C₆)alkyl-, R²³R²⁴N-(C=O)-, R²¹(C=O)-NH-, R²¹(C=O)-[N-(C₁-C₆)alkyl]-; R²¹(C=O)-NH-(C₁-C₆)alkyl-; and R²¹(C=O)-[N-(C₁-C₆)alkyl]-(C₁-C₆)alkyl-; wherein said (C₃-C₁₀)cycloalkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl(CH₂)_n-, (C₁-C₉)heterocyclic substituents may optionally be substituted on a ring carbon or nitrogen by one to three members per ring independently selected from halo, (C₁-C₆)alkyl, and (C₁-C₆)alkoxy.

50. (Currently Amended) A compound according to claim 1 ~~claims 1-45~~, wherein R⁵ is (C₆-C₁₀)aryl-(C₁-C₆)alkyl-O-, (C₁-C₉)heteroaryl-(C₁-C₆)alkyl-O-, (C₁-C₉)heterocyclic-(C₁-C₆)alkyl-O-, wherein each of said (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl, (C₁-C₉)heterocyclic moieties of said (C₆-C₁₀)aryl-(C₁-C₆)alkyl-O-, (C₁-C₉)heteroaryl-(C₁-C₆)alkyl-O-, and (C₁-C₉)heterocyclic-(C₁-C₆)alkyl-O-, may optionally be substituted with a substituent selected from the group consisting of

(C₁-C₆)alkyl, (C₂-C₆)alkenyl, (C₂-C₆)alkynyl, (C₃-C₁₀)cycloalkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl(CH₂)_n-, (C₁-C₉)heterocyclic, halo, HO-, HO-(C=O)-, R²⁰-O-(C=O)-, R²¹-(C=O)-, R²²-CO₂-, N≡C-, R²³R²⁴N-, R²³R²⁴N-(C₁-C₆)alkyl-, R²³R²⁴N-(C=O)-, R²¹(C=O)-NH-, R²¹(C=O)-[N-(C₁-C₆)alkyl]-; R²¹(C=O)-NH-(C₁-C₆)alkyl-; and R²¹(C=O)-[N-(C₁-C₆)alkyl]-(C₁-C₆)alkyl-; wherein said (C₃-C₁₀)cycloalkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl(CH₂)_n-, (C₁-C₉)heterocyclic substituents may optionally be substituted on a ring carbon or nitrogen by one to three members per ring independently selected from halo, (C₁-C₆)alkyl, and (C₁-C₆)alkoxy.

51. (Currently Amended) A compound according to claim 1 ~~claims 1-45~~, wherein R⁵ is (C₁-C₉)heteroaryl-(C₁-C₆)alkyl-O- optionally substituted with one to two substituents independently selected from the group consisting of (C₁-C₆)alkyl, (C₂-C₆)alkenyl, (C₂-C₆)alkynyl, (C₃-C₁₀)cycloalkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl(CH₂)_n-, (C₁-C₉)heterocyclic, halo, HO-, HO-(C=O)-, R²⁰-O-(C=O)-, R²¹-(C=O)-, R²²-CO₂-, N≡C-, R²³R²⁴N-, R²³R²⁴N-(C₁-C₆)alkyl-, R²³R²⁴N-(C=O)-, R²¹(C=O)-NH-, R²¹(C=O)-[N-(C₁-C₆)alkyl]-; R²¹(C=O)-NH-(C₁-C₆)alkyl-; and R²¹(C=O)-[N-(C₁-C₆)alkyl]-(C₁-C₆)alkyl-; wherein said (C₃-C₁₀)cycloalkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl(CH₂)_n-, (C₁-C₉)heterocyclic substituents may optionally be substituted on a ring carbon or nitrogen by one to three members per ring independently selected from halo, (C₁-C₆)alkyl, and (C₁-C₆)alkoxy.

52. (Currently Amended) A compound according to claim 1 ~~claims 1-45~~, wherein R⁵ is (C₁-C₉)heteroaryl-(C₁-C₆)alkyl-O- optionally substituted with one to two substituents independently selected from the group consisting of (C₁-C₆)alkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl(CH₂)_n-, halo, HO-, HO-(C=O)-, R²⁰-O-(C=O)-, R²¹-(C=O)-, R²²-CO₂-, N≡C-, R²³R²⁴N-, R²³R²⁴N-(C₁-C₆)alkyl-, R²³R²⁴N-(C=O)-, R²¹(C=O)-NH-, R²¹(C=O)-[N-(C₁-C₆)alkyl]-; R²¹(C=O)-NH-(C₁-C₆)alkyl-; and R²¹(C=O)-[N-(C₁-C₆)alkyl]-(C₁-C₆)alkyl-; wherein said (C₃-C₁₀)cycloalkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl(CH₂)_n-, (C₁-C₉)heterocyclic substituents may optionally be substituted on a ring carbon or nitrogen by one to two members per ring independently selected from halo, (C₁-C₆)alkyl, and (C₁-C₆)alkoxy;

wherein n is an integer from zero to two;

wherein each of R²³ and R²⁴ is independently selected from hydrogen, (C₁-C₆)alkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl, (C₁-C₉)heterocyclic, (C₁-C₉)heteroaryl(C₁-C₆)alkyl, (C₆-C₁₀)aryl(C₁-C₆)alkyl, (C₁-C₉)heterocyclic(C₁-C₆)alkyl, HO-(C₁-C₆)alkyl, amino-(C₁-C₆)alkyl-, (C₁-C₆)alkylamino-(C₁-C₆)alkyl-, and [(C₁-C₆)alkyl]₂amino-(C₁-C₆)alkyl-; wherein said each of said (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl, and (C₁-C₉)heterocyclic moieties of said (C₆-C₁₀)aryl-, (C₁-C₉)heteroaryl-, (C₁-C₉)heterocyclic-, (C₆-C₁₀)aryl-(C₁-C₆)alkyl, (C₁-C₉)heteroaryl-(C₁-C₆)alkyl and (C₁-C₉)heterocyclic-(C₁-C₆)alkyl, may optionally be substituted with one to two substituents independently selected from the group consisting of halo, (C₁-C₆)alkyl or (C₁-C₆)alkoxy, or R²³ and R²⁴ are taken together to form an azetidiny, pyrrolidiny, piperidiny, piperaziny, (C₁-C₆)alkyl-piperaziny or morpholiny ring.

53. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-45~~, wherein R⁵ is optionally substituted (C₁-C₆)alkyl-O-.

54. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-45~~, wherein R⁵ is (C₁-C₆)alkyl-O- optionally substituted with one to three substituents independently selected from the group consisting of (C₃-C₁₀)cycloalkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl and (C₁-C₉)heterocyclic; wherein said (C₃-C₁₀)cycloalkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl(CH₂)_n-, (C₁-C₉)heterocyclic substituents may optionally be substituted on a ring carbon or nitrogen by one to three members per ring independently selected from halo, (C₁-C₆)alkyl, and (C₁-C₆)alkoxy.

55. (Withdrawn-Currently Amended) A compound according to claim 1 ~~claims 1-45~~, wherein R⁵ is (C₁-C₆)alkyl-O- substituted with one substituent selected from the group consisting of halo, HO-, HO-(C=O)-, R²⁰-O-(C=O)-, R²¹-(C=O)-, R²²-CO₂-, N≡C-, R²³R²⁴N-, R²³R²⁴N-(C=O)-, R²¹(C=O)-NH-, and R²¹(C=O)-[N-(C₁-C₆)alkyl]-; wherein R²³ and R²⁴ is independently selected from hydrogen, (C₁-C₆)alkyl, (C₆-C₁₀)aryl, (C₁-C₉)heteroaryl, (C₁-C₉)heterocyclic, (C₁-C₉)heteroaryl(C₁-C₆)alkyl, (C₆-C₁₀)aryl(C₁-C₆)alkyl, (C₁-C₉)heterocyclic(C₁-C₆)alkyl, HO-(C₁-C₆)alkyl,

N≡C-(C₁-C₆)alkyl-, amino-(C₁-C₆)alkyl-, (C₁-C₆)alkylamino-(C₁-C₆)alkyl-, and [(C₁-C₆)alkyl]₂amino-(C₁-C₆)alkyl-; wherein said each of said (C₆-C₁₀)aryl-, (C₁-C₉)heteroaryl-, and (C₁-C₉)heterocyclic moieties of said (C₆-C₁₀)aryl-, (C₁-C₉)heteroaryl-, (C₁-C₉)heterocyclic-, (C₆-C₁₀)aryl-(C₁-C₆)alkyl-, (C₁-C₉)heteroaryl-(C₁-C₆)alkyl and (C₁-C₉)heterocyclic-(C₁-C₆)alkyl, may optionally be substituted with one to two substituents independently selected from the group consisting of halo, (C₁-C₆)alkyl or (C₁-C₆)alkoxy, or R²³ and R²⁴ are taken together to form an azetidiny, pyrrolidiny, piperidiny or morpholinyl ring.

56. (Withdrawn-Currently Amended) A compound according to claims 1, 4, 5, 6, and 7 ~~claims 1-45~~, wherein R⁵ is -C≡N, R¹⁶R¹⁷N-(C=O)-, R¹⁶R¹⁷-N-SO₂-, R¹⁸-SO₂-, R¹⁸-SO₂-(NR¹⁹)-, R¹⁸-SO₃-, R¹⁶-(C=O)-(R²⁵-N)-, R¹⁶R¹⁷N-(C=O)-(R²⁵-N)-, R¹⁹-O-(C=O)-(R²⁵-N)-, R¹⁸-(C=O)-O-, R¹⁸-(C=O)-, R¹⁶R¹⁷N-(C=O)-O- or R¹⁹-O-(C=O)-.

57. (Withdrawn-Currently Amended) A compound according to claims 1, 4, 5, 6, and 7 ~~claims 1-45~~, wherein R⁵ is R¹⁶R¹⁷N-(C=O)-.

58. (Currently Amended) A compound according to claim 1 ~~claims 1-57~~, wherein X and Y are each hydrogen.

59. (Currently Amended) A compound according to claim 1 ~~claims 1-57~~, wherein one of X and Y is fluoro, chloro, or bromo.

60. (Currently Amended) A compound according to claim 1 ~~claims 1-57~~, wherein each of X and Y are independently selected from hydrogen, fluoro, chloro, or bromo.

61. (Currently Amended) A compound according to claim 1 ~~claims 1-57~~, wherein one of X and Y is (C₁-C₆)alkyl.

62. (Original) A compound according to claim 1, wherein said compound is

(2R, 3S, 4aR, 10aR)-4a-Ethyl-2-prop-1-ynyl-1,2,3,4,4a,9,10,10a-octahydrophenanthrene-2,3,7-triol;

(2R, 3S, 4aR, 10aR)-4a-Ethyl-7-(2-methylpyridin-3-ylmethoxy)-2-prop-1-ynyl-1,2,3,4,4a,9,10,10a-octahydrophenanthrene-2,3-diol;

(2R, 3R, 4aR, 10aR)-7-[5-(2-Dimethylaminoethyl)-[1,2,4]oxadiazol-3-ylmethoxy]-4a-ethyl-3-methyl-2-phenyl-1,2,3,4,4a,9,10,10a-octahydrophenanthrene-2,3-diol

(2R, 3R, 4aR, 10aR)-4a-Ethyl-3-methyl-2-pyridin-2-yl-1,2,3,4,4a,9,10,10a-octahydrophenanthrene-2,3,7-triol;

(2R, 3R, 4aR, 10aR)-4a-Ethyl-3-methyl-7-(2-methylpyridin-3-ylmethoxy)-2-pyridin-2-yl-1,2,3,4,4a,9,10,10a-octahydrophenanthrene-2,3-diol;

(2R, 3S, 4aR, 10aR)-4a-Ethyl-3-methyl-2-thiazol-2-yl-1,2,3,4,4a,9,10,10a-octahydrophenanthrene-2,3,7-triol;

(2R, 3S, 4aR, 10aR)-4a-Ethyl-3-methyl-2-(4-methylthiazol-2-yl)-1,2,3,4,4a,9,10,10a-octahydrophenanthrene-2,3,7-triol;

(2R, 3R, 4aR, 10aS)-4a-Ethyl-2,3,7-trihydroxy-3-methyl-2-phenyl-2,3,4,4a,10,10a-hexahydro-1H-phenanthren-9-one;

(2R, 3R, 4aR, 10aS)-4a-Ethyl-3,9-dimethyl-2-phenyl-1,2,3,4,4a,10a-hexahydro-phenanthrene-2,3,7-triol;

(2R, 3R, 4aR, 10aR)-3,4a-Diethyl-2-phenyl-1,2,3,4,4a,9,10,10a-octahydro-phenanthrene-2,3,7-triol;

(2R, 3R, 4aR, 10aR)-4a-Ethyl-7-(2-hydroxy-ethoxy)-3-methyl-2-phenyl-1,2,3,4,4a,9,10,10a-octahydro-phenanthrene-2,3-diol;

(2R, 3R, 4aR, 10aR)-4a-Ethyl-7-(3-hydroxy-propoxy)-3-methyl-2-phenyl-1,2,3,4,4a,9,10,10a-octahydro-phenanthrene-2,3-diol;

(2R, 3R, 4aR, 10aR)-4a-Ethyl-7-(4-hydroxy-butoxy)-3-methyl-2-phenyl-1,2,3,4,4a,9,10,10a-octahydro-phenanthrene-2,3-diol;

(4bR, 7R, 6R, 8aR)-4-(4b-Ethyl-6,7-dihydroxy-6-methyl-7-phenyl-4b,5,6,7,8,8a,9,10-octahydro-phenanthren-2-yloxy)-butyronitrile;

(4bR, 7R, 6R, 8aR)-5-(4b-Ethyl-6,7-dihydroxy-6-methyl-7-phenyl-4b,5,6,7,8,8a,9,10-octahydro-phenanthren-2-yloxy)-pentanenitrile;

(4bR, 7R, 6R, 8aR)-2-(4*b*-Ethyl-6,7-dihydroxy-6-methyl-7-phenyl-4*b*,5,6,7,8,8*a*,9,10-octahydro-phenanthren-2-yloxy)-acetamide;
(2R, 3R, 4aR, 10aR)-4*a*-Ethyl-7-(4-hydroxy-4-methyl-pentyloxy)-3-methyl-2-phenyl-1,2,3,4,4*a*,9,10,10*a*-octahydro-phenanthrene-2,3-diol;
(2R, 3R, 4aR, 10aR)-4*a*-Ethyl-7-(5-hydroxy-5-methyl-hexyloxy)-3-methyl-2-phenyl-1,2,3,4,4*a*,9,10,10*a*-octahydro-phenanthrene-2,3-diol;
(2R, 3R, 4aR, 10aR)-4*a*-Ethyl-3-methyl-2-prop-1-ynyl-1,2,3,4,4*a*,9,10,10*a*-octahydro-phenanthrene-2,3,7-triol;
(2R, 3R, 4aR, 10aR)-4*a*-Ethyl-3-methyl-2-p-tolyl-1,2,3,4,4*a*,9,10,10*a*-octahydro-phenanthrene-2,3,7-triol; and
(2R, 3R, 4aR, 10aR)-4*a*-Ethyl-3-methyl-2-propenyl-1,2,3,4,4*a*,9,10,10*a*-octahydro-phenanthrene-2,3,7-triol.

63. (Withdrawn) A method of treating a disorder selected from the group consisting of inflammatory disorders, endocrine disorders; collagen diseases; dermatologic diseases; allergic states; ophthalmic diseases; respiratory diseases; hematologic disorders; neoplastic diseases; edematous states; and gastrointestinal diseases in a mammal comprising administering to said mammal a therapeutically effective amount of a compound according to claim 1.

64. (Original) A pharmaceutical composition for treating a disorder selected from the group consisting of inflammatory disorders, endocrine disorders; collagen diseases; dermatologic diseases; allergic states; ophthalmic diseases; respiratory diseases; hematologic disorders; neoplastic diseases; edematous states; and gastrointestinal diseases in a mammal comprising a therapeutically effective amount of a compound according to claim 1 or a salt or prodrug thereof, and a pharmaceutically acceptable carrier.

65. (Withdrawn) A method of treating inflammation in a mammal comprising administering to said mammal a therapeutically effective amount of a compound of

claim 1, an isomer thereof, a prodrug of said compound or isomer, or a pharmaceutically acceptable salt of said compound, isomer or prodrug.

66. (Original) A pharmaceutical composition for the treatment of inflammation comprising an amount of a compound of claim 1 effective for treating inflammation, an isomer thereof, a prodrug of said compound or isomer, or a pharmaceutically acceptable salt of said compound, isomer or prodrug; and a pharmaceutically acceptable carrier, vehicle or diluent.

67. (New) A compound according to claims 1, 4, 5, 6, and 7, wherein

R⁴ is HO-; and

R⁵ is -C≡N, R¹⁶R¹⁷N-(C=O)-, R¹⁶R¹⁷-N-SO₂-, R¹⁸-SO₂-, R¹⁸-SO₂-(NR¹⁹)-, R¹⁸-SO₃-, R¹⁶-(C=O)-(R²⁵-N)-, R¹⁶R¹⁷N-(C=O)-(R²⁵-N)-, R¹⁹-O-(C=O)-(R²⁵-N)-, R¹⁸-(C=O)-O-, R¹⁸-(C=O)-, R¹⁶R¹⁷N-(C=O)-O- or R¹⁹-O-(C=O)-.